

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120219\
 Data File : BN008813.D
 Acq On : 02 Dec 2019 14:19
 Operator : JU
 Sample : K5851-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AA8

Quant Time: Dec 02 15:11:18 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 02 12:46:36 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	5298	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	21588	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	12711	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	25064	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	17031	0.40	ng/ul	0.01
20) Perylene-d12	23.32	264	18195	0.40	ng/ul	0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.94	152	4670	0.13	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	10383	0.14	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.39	128	89339	1.373	ng/ul	99
5) 2-Methylnaphthalene	12.01	142	39886	0.878	ng/ul	100
7) Acenaphthylene	13.93	152	337973	4.809	ng/ul	100
8) Acenaphthene	14.27	153	80199	1.538	ng/ul	99
9) Fluorene	15.27	166	520091	8.493	ng/ul	99
12) Phenanthrene	17.00	178	4634753	54.831	ng/ul	100
13) Anthracene	17.09	178	1184030	14.825	ng/ul	99
15) Fluoranthene	19.03	202	6416702	61.789	ng/ul	96
17) Pyrene	19.39	202	5097755	74.610	ng/ul	99
18) Benzo(a)anthracene	21.15	228	2256168	32.941	ng/ul	96
19) Chrysene	21.20	228	2097239	31.176	ng/ul	96
21) Benzo(b)fluoranthene	22.69	252	3025803	40.328	ng/ul	95
22) Benzo(k)fluoranthene	22.85	252	343522	4.616	ng/ul	96
23) Benzo(a)pyrene	23.14	252	778663	11.057	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.46	276	892812	10.840	ng/ul	90
25) Dibenzo(a,h)anthracene	25.47	278	219867	3.305	ng/ul	96
26) Benzo(g,h,i)perylene	26.11	276	846874	12.292	ng/ul	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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